

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062423\
 Data File : BN025989.D
 Acq On : 24 Jun 2023 04:10
 Operator : MA/JU
 Sample : PB153403BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS403

Quant Time: Jun 24 04:49:02 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 01:27:41 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.980	152	6557	0.400	ng/ul	0.00
4) Naphthalene-d8	10.792	136	21153	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.617	164	12958	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.365	188	27366	0.400	ng/ul	0.00
17) Chrysene-d12	21.544	240	18613	0.400	ng/ul	0.00
23) Perylene-d12	24.014	264	15368	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	4909	0.665	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.376	152	12407	0.381	ng/ul	0.00
18) Fluoranthene-d10	19.387	212	28745	0.406	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.364	88	11996	1.530	ng/ul#	86
5) Naphthalene	10.836	128	21222	0.350	ng/ul	99
7) 2-Methylnaphthalene	12.453	142	13361	0.346	ng/ul	100
8) 1-Methylnaphthalene	12.667	142	14202	0.355	ng/ul	100
10) Acenaphthylene	14.340	152	22167	0.365	ng/ul	100
11) Acenaphthene	14.677	153	17255	0.348	ng/ul	98
12) Fluorene	15.663	166	19219	0.338	ng/ul	99
14) Pentachlorophenol	17.014	266	5317	0.542	ng/ul	99
15) Phenanthrene	17.403	178	30884	0.342	ng/ul	99
16) Anthracene	17.495	178	27977	0.348	ng/ul	99
19) Fluoranthene	19.415	202	36061	0.363	ng/ul	99
20) Pyrene	19.778	202	37514	0.363	ng/ul	99
21) Benzo(a)anthracene	21.530	228	28001	0.365	ng/ul	99
22) Chrysene	21.579	228	28606	0.358	ng/ul	99
24) Benzo(b)fluoranthene	23.257	252	23556	0.359	ng/ul	97
25) Benzo(k)fluoranthene	23.304	252	24791	0.376	ng/ul	98
26) Benzo(a)pyrene	23.906	252	20481	0.360	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.605	276	24845	0.399	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.615	278	19043	0.404	ng/ul	97
29) Benzo(g,h,i)perylene	27.393	276	20979	0.379	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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