

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101623\
 Data File : BN028294.D
 Acq On : 16 Oct 2023 17:26
 Operator : MA/JU
 Sample : PB156403BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS403

Quant Time: Oct 17 02:10:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 16:16:04 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.895	152	2535	0.400	ng/ul	0.00
4) Naphthalene-d8	10.713	136	8557	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.541	164	5195	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.303	188	11560	0.400	ng/ul	0.00
17) Chrysene-d12	21.501	240	11099	0.400	ng/ul	0.00
23) Perylene-d12	23.953	264	12018	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.262	96	2115	0.664	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.308	152	4440	0.322	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	11340	0.322	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.296	88	6710	1.990	ng/ul#	87
5) Naphthalene	10.768	128	7912	0.326	ng/ul	97
7) 2-Methylnaphthalene	12.379	142	5046	0.329	ng/ul	99
8) 1-Methylnaphthalene	12.594	142	4986	0.314	ng/ul	100
10) Acenaphthylene	14.273	152	7839	0.333	ng/ul	99
11) Acenaphthene	14.606	153	6194	0.330	ng/ul	100
12) Fluorene	15.600	166	7074	0.339	ng/ul	99
14) Pentachlorophenol	16.953	266	1420	0.573	ng/ul	100
15) Phenanthrene	17.346	178	11676	0.339	ng/ul	98
16) Anthracene	17.447	178	10076	0.327	ng/ul	99
19) Fluoranthene	19.366	202	13948	0.329	ng/ul	99
20) Pyrene	19.734	202	14530	0.331	ng/ul	99
21) Benzo(a)anthracene	21.484	228	14016	0.346	ng/ul	99
22) Chrysene	21.539	228	14858	0.363	ng/ul	100
24) Benzo(b)fluoranthene	23.190	252	15249	0.342	ng/ul	99
25) Benzo(k)fluoranthene	23.243	252	15943	0.360	ng/ul	99
26) Benzo(a)pyrene	23.845	252	14303	0.376	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.532	276	17777	0.387	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.545	278	14476	0.387	ng/ul	99
29) Benzo(g,h,i)perylene	27.334	276	14734	0.392	ng/ul	98

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

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