

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091523\
 Data File : BN027723.D
 Acq On : 16 Sep 2023 09:02
 Operator : MA/JU
 Sample : PB155391BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS391

Quant Time: Sep 18 00:49:52 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 16 01:11:24 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.013	152	3894	0.400	ng/ul	0.00
4) Naphthalene-d8	10.834	136	12209	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.647	164	6866	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.401	188	16972	0.400	ng/ul	0.00
17) Chrysene-d12	21.580	240	14745	0.400	ng/ul	0.00
23) Perylene-d12	24.073	264	13231	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	4093	0.834	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.412	152	8132	0.440	ng/ul	0.00
18) Fluoranthene-d10	19.422	212	21235	0.472	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.431	88	10239	2.006	ng/ul#	84
5) Naphthalene	10.883	128	14202	0.407	ng/ul	100
7) 2-Methylnaphthalene	12.483	142	8854	0.400	ng/ul	100
8) 1-Methylnaphthalene	12.703	142	9640	0.422	ng/ul	99
10) Acenaphthylene	14.379	152	12444	0.415	ng/ul	99
11) Acenaphthene	14.712	153	10863	0.419	ng/ul	100
12) Fluorene	15.697	166	12727	0.417	ng/ul	99
14) Pentachlorophenol	17.029	266	2785	0.701	ng/ul	99
15) Phenanthrene	17.443	178	22094	0.414	ng/ul	99
16) Anthracene	17.536	178	18227	0.408	ng/ul	99
19) Fluoranthene	19.450	202	25570	0.429	ng/ul	99
20) Pyrene	19.817	202	25831	0.421	ng/ul	99
21) Benzo(a)anthracene	21.559	228	22827	0.400	ng/ul	100
22) Chrysene	21.615	228	24376	0.413	ng/ul	99
24) Benzo(b)fluoranthene	23.296	252	24176	0.410	ng/ul	99
25) Benzo(k)fluoranthene	23.348	252	23616	0.413	ng/ul	99
26) Benzo(a)pyrene	23.959	252	19232	0.405	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.706	276	24066	0.396	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.727	278	19354	0.395	ng/ul	99
29) Benzo(g,h,i)perylene	27.525	276	20680	0.401	ng/ul	99

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

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