

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
 Data File : BP013339.D
 Acq On : 27 Dec 2022 15:26
 Operator : CG/JU
 Sample : N6133-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXGJ1

Quant Time: Dec 28 00:19:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 28 00:16:35 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	509247	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	2185809	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	1419403	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	3058103	20.000	ng/u1	0.00
79) Chrysene-d12	21.422	240	2618247	20.000	ng/u1	0.00
88) Perylene-d12	23.892	264	2789354	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	48594	4.080	ng/uL	0.00
4) Pyridine-d5	3.752	84	373381	11.036	ng/u1	0.00
7) Phenol-d5	7.093	99	258790	6.239	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	727933	29.092	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	721142	22.085	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	405060	11.918	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	464067	28.896	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	517569	28.626	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	963494	27.436	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	420653	8.485	ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	3492612	32.017	ng/u1	0.00
49) Acenaphthylene-d8	14.269	160	3589685	29.947	ng/u1	0.00
54) 4-Nitrophenol-d4	14.787	143	100778	5.136	ng/u1	0.00
60) Fluorene-d10	15.569	176	2938233	31.837	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	577263	29.333	ng/u1	0.00
73) Anthracene-d10	17.434	188	4606014	33.385	ng/u1	0.00
81) Pyrene-d10	19.669	212	5237855	34.852	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	4595393	33.053	ng/u1	0.00

Target Compounds Qvalue

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

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