

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
 Data File : BP013342.D
 Acq On : 27 Dec 2022 17:09
 Operator : CG/JU
 Sample : AR1254 50ppm
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 28 00:20:13 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	7.928	152	541701	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	2012653	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	1271880	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.333	188	3013998	20.000	ng/u1	0.00
79) Chrysene-d12	21.421	240	2047855	20.000	ng/u1	0.00
88) Perylene-d12	23.886	264	1943801	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	91	0.007	ng/uL	-0.09
4) Pyridine-d5	3.740	84	341	0.009	ng/u1	0.00
7) Phenol-d5	7.075	99	19	0.000	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.257	67	158	0.006	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	34	0.001	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	31	0.001	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	131	0.009	ng/u1	0.02
24) 2-Nitrophenol-d4	9.822	143	154	0.009	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	86	0.003	ng/u1	0.08
31) 4-Chloroaniline-d4	10.893	131	107	0.002	ng/u1	0.00
46) Dimethylphthalate-d6	13.992	166	70	0.001	ng/u1	0.01
49) Acenaphthylene-d8	14.269	160	142	0.001	ng/u1	0.00
54) 4-Nitrophenol-d4	14.781	143	54	0.003	ng/u1	0.00
60) Fluorene-d10	15.581	176	1094	0.013	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.704	200	37	0.002	ng/u1	0.01
73) Anthracene-d10	17.333	188	3013998	22.166	ng/u1	-0.10
81) Pyrene-d10	19.663	212	7992	0.068	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.886	264	1742706	17.987	ng/u1	0.15

Target Compounds Qvalue

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

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