

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
 Data File : BP013341.D
 Acq On : 27 Dec 2022 16:35
 Operator : CG/JU
 Sample : AR1248 50ppm
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 28 00:19:58 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	535836	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	1981207	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	1226458	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.333	188	2825742	20.000	ng/u1	0.00
79) Chrysene-d12	21.421	240	2369026	20.000	ng/u1	0.00
88) Perylene-d12	23.886	264	2250326	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	213	0.017	ng/uL	0.00
4) Pyridine-d5	3.752	84	91	0.003	ng/u1	0.00
7) Phenol-d5	7.081	99	27	0.001	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.252	67	541	0.021	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	89	0.003	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	71	0.002	ng/u1	-0.01
21) Nitrobenzene-d5	9.234	128	167	0.011	ng/u1	0.14
24) 2-Nitrophenol-d4	9.840	143	13	0.001	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.446	165	54	0.002	ng/u1	0.08
31) 4-Chloroaniline-d4	10.869	131	83	0.002	ng/u1	-0.02
46) Dimethylphthalate-d6	13.981	166	76	0.001	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	207	0.002	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.787	143	9	0.001	ng/u1	0.00
60) Fluorene-d10	15.569	176	1303	0.016	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	25	0.001	ng/u1	-0.01
73) Anthracene-d10	17.333	188	2825742	22.166	ng/u1	-0.10
81) Pyrene-d10	19.663	212	9688	0.071	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.886	264	1995344	17.789	ng/u1	0.15

Target Compounds Qvalue

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

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