

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021025\
 Data File : BP024020.D
 Acq On : 10 Feb 2025 20:34
 Operator : RC/JU
 Sample : PB166622BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK622

Quant Time: Feb 12 15:01:51 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 11:19:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	383711	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.534	136	1595931	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.381	164	963932	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	2013958	20.000	ng/u1	0.00
79) Chrysene-d12	21.604	240	1905760	20.000	ng/u1	-0.04
88) Perylene-d12	24.945	264	1921257	20.000	ng/u1	-0.06
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	61043	6.610	ng/uL	0.00
4) Pyridine-d5	3.699	84	874018	32.798	ng/u1	0.00
7) Phenol-d5	6.946	99	1073240	31.606	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.099	67	592833	32.983	ng/u1	0.00
11) 2-Chlorophenol-d4	7.310	132	867356	32.527	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	836648	30.161	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	411251	32.208	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	414916	30.116	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.163	165	768578	30.409	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	1146566	30.333	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	2264988	31.502	ng/u1	-0.01
49) Acenaphthylene-d8	14.069	160	2711905	33.442	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.604	143	351951	24.648	ng/u1	0.00
60) Fluorene-d10	15.386	176	2052641	33.901	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	215270	17.326	ng/u1	0.00
73) Anthracene-d10	17.280	188	3427469	34.661	ng/u1	0.00
81) Pyrene-d10	19.645	212	3857934	37.781	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.715	264	3525473	35.169	ng/u1	-0.06

Target Compounds Qvalue

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

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