

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021025\
 Data File : BP024008.D
 Acq On : 10 Feb 2025 11:48
 Operator : RC/JU
 Sample : PB166626BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK626

Quant Time: Feb 12 14:27:00 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	7.746	152	365008	20.000	ng/u1	-0.03
20) Naphthalene-d8	10.528	136	1638732	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.381	164	1044251	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.175	188	2014118	20.000	ng/u1	-0.01
79) Chrysene-d12	21.610	240	1721484	20.000	ng/u1	-0.03
88) Perylene-d12	24.974	264	1642884	20.000	ng/u1	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	59482	6.771	ng/uL	-0.02
4) Pyridine-d5	3.687	84	885969	34.950	ng/u1	-0.02
7) Phenol-d5	6.934	99	1186556	36.733	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.081	67	597179	34.927	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.293	132	947271	37.344	ng/u1	-0.02
15) 4-Methylphenol-d8	8.452	113	935567	35.455	ng/u1	-0.02
21) Nitrobenzene-d5	8.899	128	433231	33.044	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.610	143	437756	30.943	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.157	165	862160	33.221	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.657	131	1255270	32.342	ng/u1	-0.02
46) Dimethylphthalate-d6	13.787	166	2472530	31.743	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2973309	33.845	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.604	143	296090	19.141	ng/u1	0.00
60) Fluorene-d10	15.386	176	2241738	34.176	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	92227	7.422	ng/u1	0.00
73) Anthracene-d10	17.269	188	3395726	34.337	ng/u1	-0.02
81) Pyrene-d10	19.639	212	3671658	39.805	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.733	264	3090688	36.056	ng/u1	-0.04

Target Compounds Qvalue

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

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