

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021633.D
 Acq On : 26 Aug 2024 10:11
 Operator : RC/JU
 Sample : PB162928BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK928

Quant Time: Aug 26 23:51:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 23 17:51:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	282406	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1248208	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	840393	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.228	188	1858453	20.000	ng/u1	-0.01
79) Chrysene-d12	21.692	240	1837381	20.000	ng/u1	-0.02
88) Perylene-d12	25.104	264	1885702	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	49699	5.843	ng/uL	0.00
4) Pyridine-d5	3.693	84	663661	26.974	ng/u1	0.00
7) Phenol-d5	6.958	99	824292	27.350	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	480693	29.308	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	531041	27.292	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	629767	27.041	ng/u1	-0.01
21) Nitrobenzene-d5	8.940	128	269297	26.816	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	268268	26.317	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	502781	25.029	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	731939	25.045	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	1831462	28.481	ng/u1	-0.01
49) Acenaphthylene-d8	14.122	160	1991776	27.588	ng/u1	0.00
54) 4-Nitrophenol-d4	14.628	143	289545	23.667	ng/u1	0.00
60) Fluorene-d10	15.439	176	1509539	27.922	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	211213	20.663	ng/u1	-0.01
73) Anthracene-d10	17.334	188	2460907	27.901	ng/u1	-0.01
81) Pyrene-d10	19.704	212	2995294	28.518	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.868	264	3082117	30.652	ng/u1	-0.04

Target Compounds	Qvalue

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

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