

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016814.D
 Acq On : 29 Aug 2023 04:19
 Operator : MA/JU
 Sample : PB155098BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK098

Quant Time: Aug 29 05:13:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 29 01:12:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	320976	20.000	ng/u1	0.00
20) Naphthalene-d8	10.857	136	1455522	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.681	164	916770	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.451	188	2044512	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	1685314	20.000	ng/u1	0.00
88) Perylene-d12	24.116	264	1526487	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	76337	9.829	ng/uL	0.00
4) Pyridine-d5	3.817	84	920705	38.794	ng/u1	0.00
7) Phenol-d5	7.169	99	1099238	38.514	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	687765	37.809	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	821988	38.304	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	873871	37.506	ng/u1	0.00
21) Nitrobenzene-d5	9.228	128	426149	38.723	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	457483	39.437	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	814073	37.643	ng/u1	0.00
31) 4-Chloroaniline-d4	11.016	131	1209647	36.474	ng/u1	0.00
46) Dimethylphthalate-d6	14.093	166	2648235	38.304	ng/u1	0.00
49) Acenaphthylene-d8	14.381	160	3075705	39.353	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	433383	30.328	ng/u1	0.00
60) Fluorene-d10	15.675	176	2330124	40.022	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	323564	26.575	ng/u1	0.00
73) Anthracene-d10	17.551	188	3780470	41.716	ng/u1	0.00
81) Pyrene-d10	19.780	212	4288774	46.335	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.945	264	3196025	41.798	ng/u1	0.00

Target Compounds Qvalue

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

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