

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014492.D
 Acq On : 03 Apr 2023 18:05
 Operator : CG/JU
 Sample : PB151858BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK858

Quant Time: Apr 04 01:38:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 04 01:34:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	114055	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	513944	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.663	164	352894	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	834705	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	791477	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	760666	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	16144	5.489	ng/uL	0.00
4) Pyridine-d5	3.816	84	218970	29.052	ng/ul	0.00
7) Phenol-d5	7.175	99	298657	28.393	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	193325	28.646	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	223553	29.405	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	245856	28.976	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	112109	27.021	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	131157	27.607	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	229892	26.570	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	321175	28.181	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	821694	28.717	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	903600	28.460	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	96168	22.247	ng/ul	0.01
60) Fluorene-d10	15.657	176	708958	29.712	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	116019	18.890	ng/ul	0.00
73) Anthracene-d10	17.522	188	1175959	29.111	ng/ul	0.00
81) Pyrene-d10	19.751	212	1429438	26.888	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.862	264	1186784	29.366	ng/ul	0.00

Target Compounds Qvalue

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

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