

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090723\
 Data File : BP016992.D
 Acq On : 07 Sep 2023 17:35
 Operator : MA/JU
 Sample : PB155068BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK068

Quant Time: Sep 08 01:22:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	278947	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1229761	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	759102	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	1627700	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	1294539	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	1289251	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	64893	8.800	ng/uL	0.00
4) Pyridine-d5	3.758	84	773035	36.686	ng/u1	0.00
7) Phenol-d5	7.128	99	895259	35.130	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	592409	37.163	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	671665	36.416	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	702661	36.192	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	339776	35.614	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	369146	35.693	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	632610	33.132	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	969238	35.185	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	2141849	37.905	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	2348378	36.012	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	373818	33.868	ng/u1	0.00
60) Fluorene-d10	15.633	176	1808431	37.380	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	261848	26.732	ng/u1	0.00
73) Anthracene-d10	17.504	188	2862309	38.291	ng/u1	0.00
81) Pyrene-d10	19.745	212	3148853	39.935	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.880	264	2592916	39.643	ng/u1	0.00

Target Compounds Qvalue

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

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