

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
 Data File : BP015413.D
 Acq On : 07 Jun 2023 23:10
 Operator : MA/JU
 Sample : PB153099BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK099

Quant Time: Jun 08 01:28:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 07 23:46:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.099	152	119324	20.000	ng/u1	0.00
20) Naphthalene-d8	10.928	136	578901	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.740	164	440211	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.498	188	1067410	20.000	ng/u1	0.00
79) Chrysene-d12	21.574	240	1135633	20.000	ng/u1	-0.01
88) Perylene-d12	24.133	264	1216126	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	22466	6.972	ng/uL	0.00
4) Pyridine-d5	3.852	84	283888	33.917	ng/u1	0.00
7) Phenol-d5	7.252	99	374436	34.062	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	252311	38.432	ng/u1	0.00
11) 2-Chlorophenol-d4	7.622	132	304956	38.262	ng/u1	0.00
15) 4-Methylphenol-d8	8.810	113	328202	36.378	ng/u1	0.00
21) Nitrobenzene-d5	9.275	128	166250	36.579	ng/u1	0.00
24) 2-Nitrophenol-d4	10.004	143	188952	36.405	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.546	165	314412	33.208	ng/u1	0.00
31) 4-Chloroaniline-d4	11.063	131	538695	39.639	ng/u1	0.00
46) Dimethylphthalate-d6	14.145	166	1255911	36.180	ng/u1	0.00
49) Acenaphthylene-d8	14.434	160	1318115	34.726	ng/u1	0.00
54) 4-Nitrophenol-d4	14.940	143	179231	28.832	ng/u1	0.00
60) Fluorene-d10	15.734	176	1018813	34.804	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	204730	30.276	ng/u1	0.00
73) Anthracene-d10	17.598	188	1710427	35.219	ng/u1	0.00
81) Pyrene-d10	19.822	212	2097861	34.515	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.963	264	2134387	34.978	ng/u1	-0.01

Target Compounds Qvalue

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

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