

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017492.D
 Acq On : 03 Oct 2023 00:14
 Operator : MA/JU
 Sample : PB155900BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK900

Quant Time: Oct 03 00:51:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	146716	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	577243	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	335563	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.422	188	701475	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	671263	20.000	ng/u1	0.00
88) Perylene-d12	24.121	264	740802	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	25032	7.008	ng/uL	0.00
4) Pyridine-d5	3.728	84	348650	34.914	ng/u1	0.00
7) Phenol-d5	7.099	99	440522	36.559	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	276718	38.053	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	359261	37.591	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	347091	37.050	ng/u1	0.00
21) Nitrobenzene-d5	9.152	128	172303	41.289	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	189609	41.845	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	328293	38.348	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	461684	37.229	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	962282	40.031	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	1081130	39.104	ng/u1	0.00
54) 4-Nitrophenol-d4	14.875	143	113996	27.220	ng/u1	0.00
60) Fluorene-d10	15.639	176	809299	39.106	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	108602	27.469	ng/u1	0.00
73) Anthracene-d10	17.522	188	1287752	41.392	ng/u1	0.00
81) Pyrene-d10	19.774	212	1548285	42.652	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.951	264	1461289	40.970	ng/u1	0.00

Target Compounds Qvalue

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

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