

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
 Data File : BP011032.D
 Acq On : 19 Jul 2022 14:58
 Operator : CG/JU
 Sample : PB146298BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK298

Quant Time: Jul 19 23:40:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	299524	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	1195174	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.339	164	658368	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	1308938	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.233	240	1287089	20.000	ng/u1	# 0.00
88) Perylene-d12	23.586	264	1302842	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	63415	7.633	ng/uL	0.00
4) Pyridine-d5	3.587	84	765554	35.649	ng/u1	0.00
7) Phenol-d5	6.857	99	866125	35.227	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	578559	36.594	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	729730	36.731	ng/u1	0.00
15) 4-Methylphenol-d8	8.392	113	677469	34.418	ng/u1	0.00
21) Nitrobenzene-d5	8.839	128	332291	38.151	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	318379	36.942	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.098	165	558893	35.460	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	914913	35.211	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	1756251	39.030	ng/u1	0.00
49) Acenaphthylene-d8	14.033	160	2152035	40.535	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	245578	26.966	ng/u1	0.00
60) Fluorene-d10	15.339	176	1470536	38.448	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	174797	26.987	ng/u1	0.00
73) Anthracene-d10	17.210	188	2223759	39.963	ng/u1	0.00
81) Pyrene-d10	19.462	212	2553424	37.774	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.433	264	2499814	39.949	ng/u1	0.00

Target Compounds	Qvalue

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

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